

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF063018\
 Data File : BF107027.D
 Acq On : 30 Jun 2018 20:51
 Operator : JU/SJ
 Sample : J3638-03 2X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P-3

Manual Integrations
 APPROVED

Sohil
 7/5/2018 4:32:20 PM

Quant Time: Jul 02 07:21:26 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 29 14:38:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	61129	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	233205	20.00	ng	-0.01
38) Acenaphthene-d10	10.00	164	114110	20.00	ng	-0.01
63) Phenanthrene-d10	11.49	188	226679	20.00	ng	-0.01
75) Chrysene-d12	14.14	240	209850	20.00	ng	-0.02
86) Perylene-d12	15.68	264	177116	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	121731	33.72	ng	-0.01
7) Phenol-d6	6.60	99	155898	34.58	ng	-0.02
23) Nitrobenzene-d5	7.51	82	86033	20.50	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	47911	36.23	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	203323	18.64	ng	-0.02
78) Terphenyl-d14	13.07	244	258506	29.84	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.70	163	21968	2.567	ng	Qvalue # 95
70) Phenanthrene	11.51	178	60451	5.529	ng	97
74) Fluoranthene	12.71	202	149052	12.369	ng	94
77) Pyrene	12.94	202	147589	10.665	ng	98
80) Benzo(a)anthracene	14.13	228	126617	9.900	ng	99
82) Chrysene	14.17	228	120348	10.228	ng	98
85) Indeno(1,2,3-cd)pyrene	17.26	276	96854	9.700	ng	# 86
87) Benzo(b)fluoranthene	15.22	252	212392m	19.304	ng	
88) Benzo(k)fluoranthene	15.25	252	66559m	6.353	ng	
89) Benzo(a)pyrene	15.61	252	136641	13.970	ng	96
91) Benzo(g,h,i)perylene	17.74	276	95462	11.783	ng	# 86

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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